

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	115	562/586	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:21
L2	1572	((562/586) or (562/856) or (568/486) or (568/488) or (568/489) or (568/490) or (568/615) or (568/622)). CCLS.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	OFF	2006/08/31 08:45
L3	8204727	reduc\$5 or hydrogen\$6	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:24
L4	964	hydrofluoroether	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:25
L5	1940726	platinum or pt!	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:45
L6	232	(I2 or I4) and I3 and I5	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:26
L7	81722	catalyst near10 I5	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 07:28
L8	81	I6 and I7	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:40
L9	660	Picozzi.in. or Meo.in. or Tonelli.in.	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:41

EAST Search History

L10	7	I3 and I5 and I7 and I9	US-PGPUB; USPAT; USOCR; EPO; JPO; DERWENT	OR	ON	2006/08/31 08:45
L11	1454	((562/586) or (562/856) or (568/486) or (568/488) or (568/489) or (568/490) or (568/615) or (568/622)). CCLS.	US-PGPUB; USPAT; USOCR	OR	OFF	2006/08/31 08:45
L12	35047	platinum.clm.	US-PGPUB; USPAT; USOCR	OR	ON	2006/08/31 08:46
L13	16	I11 and I12	US-PGPUB; USPAT; USOCR	OR	ON	2006/08/31 08:46

10/630,697

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1204rxw

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS	1		Web Page URLs for STN Seminar Schedule - N. America
NEWS	2		"Ask CAS" for self-help around the clock
NEWS	3	FEB 27	New STN AnaVist pricing effective March 1, 2006
NEWS	4	APR 04	STN AnaVist \$500 visualization usage credit offered
NEWS	5	MAY 10	CA/CAPLUS enhanced with 1900-1906 U.S. patent records
NEWS	6	MAY 11	KOREAPAT updates resume
NEWS	7	MAY 19	Derwent World Patents Index to be reloaded and enhanced
NEWS	8	MAY 30	IPC 8 Rolled-up Core codes added to CA/CAPLUS and USPATFULL/USPAT2
NEWS	9	MAY 30	The F-Term thesaurus is now available in CA/CAPLUS
NEWS	10	JUN 02	The first reclassification of IPC codes now complete in INPADOC
NEWS	11	JUN 26	TULSA/TULSA2 reloaded and enhanced with new search and and display fields
NEWS	12	JUN 28	Price changes in full-text patent databases EPFULL and PCTFULL
NEWS	13	JUL 11	CHEMSAFE reloaded and enhanced
NEWS	14	JUL 14	FSTA enhanced with Japanese patents
NEWS	15	JUL 19	Coverage of Research Disclosure reinstated in DWPI
NEWS	16	AUG 09	INSPEC enhanced with 1898-1968 archive
NEWS	17	AUG 28	ADISCTI Reloaded and Enhanced
NEWS	18	AUG 30	CA(SM)/CAPLUS(SM) Austrian patent law changes
NEWS EXPRESS		JUNE 30	CURRENT WINDOWS VERSION IS V8.01b, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 JUNE 2006.
NEWS HOURS			STN Operating Hours Plus Help Desk Availability
NEWS LOGIN			Welcome Banner and News Items
NEWS IPC8			For general information regarding STN implementation of IPC 8
NEWS X25			X.25 communication option no longer available

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 07:02:32 ON 31 AUG 2006

=> file reg

COST IN U.S. DOLLARS

SINCE FILE
ENTRY

TOTAL
SESSION

10/630,697

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 07:02:59 ON 31 AUG 2006
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2006 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 30 AUG 2006 HIGHEST RN 905475-39-0
DICTIONARY FILE UPDATES: 30 AUG 2006 HIGHEST RN 905475-39-0

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 963

L1 SCREEN CREATED

=> screen 1992 OR 2016 OR 2021 OR 2026 OR 1838

L2 SCREEN CREATED

=>

Uploading C:\Documents and Settings\rkeys\My
Documents\STNEXP4\TEMPLATE\STANDARD\10630697.str

1111 — 1111

1111 — 1111

10/630,697

chain nodes :

1 2 3 4 5 6 7 9 10 11 12 13 15 16 17 18

chain bonds :

1-2 2-3 2-4 4-5 4-6 4-9 6-7 7-18 10-11 11-12 11-13 11-15 15-16 16-17

exact/norm bonds :

2-3 4-6 4-9 6-7 11-13 11-15 15-16

exact bonds :

1-2 2-4 4-5 7-18 10-11 11-12 16-17

G1:F,CF3

Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 9:CLASS 10:CLASS

11:CLASS 12:CLASS 13:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS

fragments assigned product role:

containing 10

fragments assigned reactant/reagent role:

containing 1

L3 STRUCTURE UPLOADED

=> que L3 AND L1 NOT L2

L4 QUE L3 AND L1 NOT L2

=> d

L4 HAS NO ANSWERS

L1 SCR 963

L2 SCR 1992 OR 2016 OR 2021 OR 2026 OR 1838

L3 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

L4 QUE L3 AND L1 NOT L2

=> file reaction

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.88

1.09

FILE 'CASREACT' ENTERED AT 07:03:55 ON 31 AUG 2006

USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT

COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'CHEMINFORMRX' ENTERED AT 07:03:55 ON 31 AUG 2006

COPYRIGHT (C) FIZ-CHEMIE BERLIN

FILE 'DJSMONLINE' ENTERED AT 07:03:55 ON 31 AUG 2006

COPYRIGHT (C) 2006 THE THOMSON CORPORATION

FILE 'PS' ENTERED AT 07:03:55 ON 31 AUG 2006

COPYRIGHT (C) 2006 Thieme on STN

=> s 14

SAMPLE SEARCH INITIATED 07:04:11 FILE 'CASREACT'

SCREENING COMPLETE - 8 REACTIONS TO VERIFY FROM

2 DOCUMENTS

100.0% DONE

8 VERIFIED

0 HIT RXNS

0 DOCS

10/630,697

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED VERIFICATIONS: 8 TO 329
PROJECTED ANSWERS: 0 TO 0

SAMPLE SEARCH INITIATED 07:04:13 FILE 'CHEMINFORMRX'
SCREENING COMPLETE - 3 REACTIONS TO VERIFY FROM 1 DOCUMENTS

100.0% DONE 3 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.07

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED VERIFICATIONS: 3 TO 163
PROJECTED ANSWERS: 0 TO 0

FULL SEARCH INITIATED 07:04:23 FILE 'DJSMONLINE'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.10

3 FILES SEARCHED...

FULL SEARCH INITIATED 07:04:35 FILE 'PS'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.03

L5 0 L4

=> s l4 ful

FULL SEARCH INITIATED 07:04:48 FILE 'CASREACT'
SCREENING COMPLETE - 153 REACTIONS TO VERIFY FROM 36 DOCUMENTS

100.0% DONE 153 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.01

FULL SEARCH INITIATED 07:04:49 FILE 'CHEMINFORMRX'
SCREENING COMPLETE - 16 REACTIONS TO VERIFY FROM 8 DOCUMENTS

100.0% DONE 16 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.12

FULL SEARCH INITIATED 07:05:02 FILE 'DJSMONLINE'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.02

3 FILES SEARCHED...

FULL SEARCH INITIATED 07:05:05 FILE 'PS'
SCREENING COMPLETE - 0 REACTIONS TO VERIFY FROM 0 DOCUMENTS

100.0% DONE 0 VERIFIED 0 HIT RXNS 0 DOCS
SEARCH TIME: 00.00.01

L6 0 L4

=> file stnguide

COST IN U.S. DOLLARS SINCE FILE TOTAL

10/630,697

	ENTRY	SESSION
FULL ESTIMATED COST	362.09	363.18

FILE 'STNGUIDE' ENTERED AT 07:05:18 ON 31 AUG 2006
USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT
COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY, JAPAN SCIENCE
AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Aug 25, 2006 (20060825/UP).

	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.18	363.36

FILE 'REGISTRY' ENTERED AT 07:07:06 ON 31 AUG 2006
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2006 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 30 AUG 2006 HIGHEST RN 905475-39-0
DICTIONARY FILE UPDATES: 30 AUG 2006 HIGHEST RN 905475-39-0

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH June 30, 2006

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=>Testing the current file.... screen

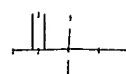
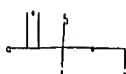
ENTER SCREEN EXPRESSION OR (END):end

=> screen 1992 OR 2016 OR 2021 OR 2026 OR 1838

L7 SCREEN CREATED

=>
Uploading C:\Documents and Settings\rkeys\My
Documents\STNEXP4\TEMPLATE\STANDARD\10630697a.str

10/630,697



chain nodes :
1 2 3 4 5 6 7 9 10
chain bonds :
1-2 2-3 2-4 4-5 4-6 4-9 6-7 7-10
exact/norm bonds :
2-3 4-6 4-9 6-7
exact bonds :
1-2 2-4 4-5 7-10

G1:F,CF3

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 9:CLASS 10:CLASS

L8 STRUCTURE UPLOADED

=> que L8 NOT L7

L9 QUE L8 NOT L7

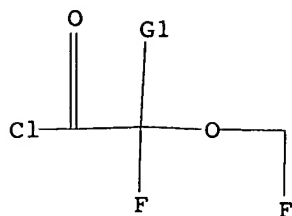
=> d

L9 HAS NO ANSWERS

L7 SCR 1992 OR 2016 OR 2021 OR 2026 OR 1838

L8 STR

10/630,697



G1 F,CF3

Structure attributes must be viewed using STN Express query preparation.
L9 QUE L8 NOT L7

=> s 19

SAMPLE SEARCH INITIATED 07:07:24 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 224 TO ITERATE

100.0% PROCESSED 224 ITERATIONS 4 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 3583 TO 5377
PROJECTED ANSWERS: 4 TO 200

L10 4 SEA SSS SAM L8 NOT L7

=> s 19 ful

FULL SEARCH INITIATED 07:07:33 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 4079 TO ITERATE

100.0% PROCESSED 4079 ITERATIONS 35 ANSWERS
SEARCH TIME: 00.00.01

L11 35 SEA SSS FUL L8 NOT L7

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	166.94	530.30

FILE 'CAPLUS' ENTERED AT 07:07:39 ON 31 AUG 2006
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 31 Aug 2006 VOL 145 ISS 10
FILE LAST UPDATED: 29 Aug 2006 (20060829/ED)

10/630,697

soluble in both fluorous solvents and liquid carbon dioxide due to the attachment of a perfluoropolyether. The enzyme horse liver alc. dehydrogenase (HLADH) was active in catalyzing oxidation/reduction reactions using FNAD as a soluble coenzyme in a fluorous solvent, methoxynonafluorobutane (HFE), and liquid carbon dioxide. In both solvents, the activity of HLADH using FNAD was greater than the same molar amount of unmodified (insol.) NAD, indicating that a soluble coenzyme results in more efficient reactions.

RE.CNT 43 THERE ARE 43 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L15 ANSWER 2 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1993:625594 CAPLUS

DN 119:225594

TI Preparation of perfluorooxaalkanoyl halides and bis(perfluorooxaalkanoyl) peroxides

IN Sawada, Hideo; Matsumoto, Takeo; Nakayama, Masaharu

PA Nippon Oils & Fats Co Ltd, Japan

SO Jpn. Kokai Tokkyo Koho, 8 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 05170731	A2	19930709	JP 1991-341515	19911224
PRAI	JP 1991-341515		19911224		
AB	RfCOX and RfCO ₂ OCORf1 [Rf, Rf1 = CF ₃ [OCF(CF ₃)CF ₂] _n (OCF ₂) _m]; X = Br, Cl, F; n, m = 1-10] are prepared Chlorination of CF ₃ OCF(CF ₃)CF ₂ OCF ₂ CO ₂ H with POCl ₃ in DMF at 100° for 5 h gave 88% CF ₃ OCF(CF ₃)CF ₂ OCF ₂ COCl, which was treated with H ₂ O ₂ in CF ₃ CF ₂ CHCl ₂ -CCl ₂ CF ₂ CHFC1 mixture at temperature between -5° and +5° to give 79% [CF ₃ OCF(CF ₃)CF ₂ OCF ₂ CO ₂] ₂ .				

L15 ANSWER 3 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1986:625978 CAPLUS

DN 105:225978

TI Asymmetric addition of hydrogen cyanide to substituted benzaldehydes catalyzed by a synthetic cyclic peptide, cyclo[(S)-phenylalanyl-(S)-histidyl]

AU Kobayashi, Yoshiyuki; Asada, Shoichi; Watanabe, Ichigen; Hayashi, Hiroaki; Motoo, Yoshiyuki; Inoue, Shohei

CS Fac. Eng., Univ. Tokyo, Tokyo, 113, Japan

SO Bulletin of the Chemical Society of Japan (1986), 59(3), 893-5

CODEN: BCSJA8; ISSN: 0009-2673

DT Journal

LA English

OS CASREACT 105:225978

AB Optically active RC₆H₄CH(OH)CN (R = H, 4-Me, 3-Me, 2-Me, 3-MeO, 3-PhO) were prepared in 33-90% enantiomeric excess by addition of HCN to RC₆H₄CHO in C₆H₆ in the presence of cyclo[(S)-phenylalanyl-(S)-histidyl]. Highest optical yields were realized in nonpolar solvents, whereas, no asym. induction occurred in MeOH or Me₂SO.

L15 ANSWER 4 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1984:103740 CAPLUS

DN 100:103740

TI Synthesis of chiral steroid CD-ring synthon from D-leucine by means of diastereotopic face selection

AU Takahashi, Takashi; Okumoto, Hiroshi; Tsuji, Jiro; Harada, Nobuyuki

CS Dep. Chem. Eng., Tokyo Inst. Technol., Tokyo, 152, Japan

SO Journal of Organic Chemistry (1984), 49(5), 948-50

CODEN: JOCEAH; ISSN: 0022-3263

- DT Journal
LA English
AB The synthesis of indanone I is described, in which the correct absolute configuration is produced from D-leucine via the cis-vinyl iodide II (R = H), which serves to control the chirality of the rest of I by means of a remarkably effective diastereotopic face-selection. The optical purity of I and II (R = H) were checked by the ^{19}F NMR of their $(\text{F}_3\text{C})\text{CFOCF}(\text{CF}_3)\text{CO}_2\text{H}$ esters. The relative and absolute configuration of I and (-)-dienone III were determined by NMR and CD data. Thus, successive treatment of II (R = MeOCMe_2) with BuLi, CuI-PBu₃, 2-methyl-2-cyclopentenone, $\text{H}_2\text{C}:\text{C}(\text{SiMe}_3)\text{COMe}$, NaOMe, and HCl gave 58% I.
- L15 ANSWER 5 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN
AN 1985:5467 CAPLUS
DN 102:5467
TI Thermal decomposition of (trifluoromethoxy)difluoroacetyl peroxide in heptane as a method of generating (trifluoromethoxy)difluoromethyl radicals
AU Komendantov, A. M.; Berenblit, V. V.; Sass, V.; Sokolov, S. V.
CS Vses. Nauchno-Issled. Inst. Sint. Kauch., USSR
SO Zhurnal Vsesoyuznogo Khimicheskogo Obshchestva im. D. I. Mendeleeva (1984), 29(3), 353-4
CODEN: ZVKOA6; ISSN: 0373-0247
DT Journal
LA Russian
AB The thermolysis of $(\text{CF}_3\text{OCF}_2\text{CO})_2\text{O}_2$ at 5-30° was a monomol., 1st-order process with activation energy 97 kJ/mol. The resulting $\text{CF}_3\text{OCF}_2\cdot$ radical abstracted H from heptane to give $\text{CF}_3\text{OCF}_2\text{H}$ quant. within 20 min at 50° with no side reaction.
- L15 ANSWER 6 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN
AN 1984:422810 CAPLUS
DN 101:22810
TI Thermal decomposition of 2-(trifluoromethoxy)perfluoropropionyl peroxide in heptane as a method for the generation of 1-(trifluoromethoxy)perfluoroethyl radicals
AU Komendantov, A. M.; Starobin, Yu. K.; Berenblit, V. V.; Sass, V. P.; Sokolov, S. V.
CS Vses. Nauchno-Issled. Inst. Sint. Kauch., Leningrad, USSR
SO Zhurnal Vsesoyuznogo Khimicheskogo Obshchestva im. D. I. Mendeleeva (1984), 29(1), 113-14
CODEN: ZVKOA6; ISSN: 0373-0247
DT Journal
LA Russian
OS CASREACT 101:22810
AB The title thermolysis at 5-25° was 1st order in peroxide. The resulting $\cdot\text{CF}(\text{CF}_3)\text{OCF}_3$ abstracted H from the solvent to give $\text{CF}_3\text{OCHFOCF}_3$.
- L15 ANSWER 7 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN
AN 1976:432486 CAPLUS
DN 85:32486
TI The reduction of perfluoroacyl halides with organosilicon hydrides. A direct synthesis of fluorine-containing esters and lactones
AU Croft, Thomas S.; McBrady, John J.
CS Cent. Res. Lab., 3M Co., St. Paul, MN, USA
SO Journal of Organic Chemistry (1976), 41(13), 2256-8
CODEN: JOCEAH; ISSN: 0022-3263
DT Journal
LA English
AB Reaction of perfluoroacyl halides with organosilicon halides in the presence of KF, ZnCl₂, and Pt/C gave 1,1-dihydroperfluoroalkyl perfluoroacylates. E.g., reaction of 2.8 g $\text{CF}_3\text{CF}_2\text{O}(\text{CF}_2)_2\text{COF}$ with Me_3SiH

10/630,697

for 18 hr at 180° in the presence of the above catalysts gave 0.5 g CF₃CF₂O(CF₂)₂CO₂CH₂(CF₂)₂OCF₂CF₃. Similarly, I was prepared from perfluorocyclohexanecarbonyl fluoride; II (x = bond, CF₂, O) were prepared from perfluorosuccinyl or perfluoroglutaryl fluoride or O(CF₂COCl)₂.

L15 ANSWER 8 OF 8 CAPLUS COPYRIGHT 2006 ACS on STN

AN 1975:594115 CAPLUS

DN 83:194115

TI Perfluorinated linear polyethers having reactive terminal groups at both ends of the chain

IN Sianesi, Dario; Caporiccio, Gerardo; Mensi, Domenico

PA Montedison S.p.A., Italy

SO U.S., 14 pp.

CODEN: USXXAM

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 3847978	A	19741112	US 1969-834486	19690618
PRAI	US 1968-787309	A2	19681226		

AB Perfluorinated linear polyethers containing peroxidic linkages were chain-cleaved by reducing agents to give bifunctional perfluorinated linear oligopolyethers with chemical-reactive terminal groups. Thus, hexafluoropropene [116-15-4] was treated with oxygen under the influence of uv light to give a peroxidized poly(perfluoropropylene oxide) [25038-02-2] which was reduced by H over a Pd catalyst to give a series of carboxy- and trifluoroacetyl-terminated oligopolyethers. One of these, CF₃COCF₂O(C₃F₆O)₂CF(CF₃)CO₂H [42775-40-6], boiling point 210-2°, formed a polymer with hexamethylenediamine [55809-69-3].

=> file stnguide

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
44.07	574.37

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-6.00	-6.00

CA SUBSCRIBER PRICE

FILE 'STNGUIDE' ENTERED AT 07:13:31 ON 31 AUG 2006

USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT

COPYRIGHT (C) 2006 AMERICAN CHEMICAL SOCIETY, JAPAN SCIENCE

AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.

LAST RELOADED: Aug 25, 2006 (20060825/UP).

=> log y

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.42	574.79

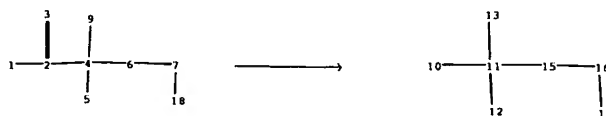
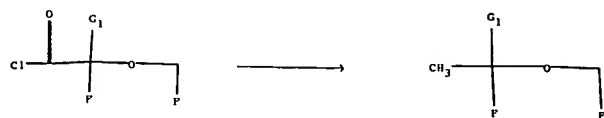
FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-6.00

CA SUBSCRIBER PRICE

STN INTERNATIONAL LOGOFF AT 07:17:27 ON 31 AUG 2006



chain nodes :

1 2 3 4 5 6 7 9 10 11 12 13 15 16 17 18

chain bonds :

1-2 2-3 2-4 4-5 4-6 4-9 6-7 7-18 10-11 11-12 11-13 11-15 15-16 16-17

exact/norm bonds :

2-3 4-6 4-9 6-7 11-13 11-15 15-16

exact bonds :

1-2 2-4 4-5 7-18 10-11 11-12 16-17

G1:F,CF3

Match level :

1:CLASS2:CLASS3:CLASS4:CLASS5:CLASS6:CLASS7:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS
15:CLASS16:CLASS17:CLASS18:CLASS

fragments assigned product role:

containing 10

fragments assigned reactant/reagent role:

containing 1